



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:39 AM GMT

PDB ID : 2B81
Title : Crystal Structure of the Luciferase-like Monooxygenase from *Bacillus cereus*
Authors : Kim, Y.; Li, H.; Moy, S.; Collart, F.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2005-10-06
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

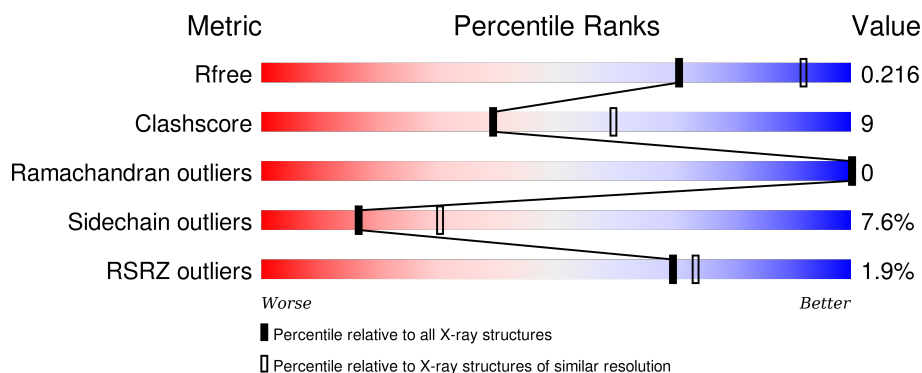
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	323	<div> <div>80%</div> <div>16%</div> <div>• •</div> </div>
1	B	323	<div> <div>82%</div> <div>13%</div> <div>• •</div> </div>
1	C	323	<div> <div>79%</div> <div>15%</div> <div>• •</div> </div>
1	D	323	<div> <div>77%</div> <div>16%</div> <div>5%</div> <div>•</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	2427	-	-	-	X
2	SO4	A	2428	-	-	-	X
2	SO4	A	2531	-	-	-	X
2	SO4	A	2532	-	-	-	X
2	SO4	D	2440	-	-	-	X
3	MES	A	2404	-	-	X	X
4	GOL	A	2415	-	-	-	X
4	GOL	B	2412	-	-	-	X
4	GOL	D	2414	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12077 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Luciferase-like monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	Se	0	14	0
			2690	1724	457	500	9			
1	B	318	Total	C	N	O	Se	0	12	0
			2690	1723	458	500	9			
1	C	320	Total	C	N	O	Se	0	9	0
			2670	1713	454	493	10			
1	D	318	Total	C	N	O	Se	0	13	0
			2698	1726	463	500	9			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	CLONING ARTIFACT	UNP Q81B18
A	-1	ASN	-	CLONING ARTIFACT	UNP Q81B18
A	0	ALA	-	CLONING ARTIFACT	UNP Q81B18
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	4	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	16	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	40	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	81	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	82	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	128	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	204	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	215	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
A	250	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	-2	SER	-	CLONING ARTIFACT	UNP Q81B18
B	-1	ASN	-	CLONING ARTIFACT	UNP Q81B18
B	0	ALA	-	CLONING ARTIFACT	UNP Q81B18
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	4	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	16	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	40	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	81	MSE	MET	MODIFIED RESIDUE	UNP Q81B18

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Chain	Residue	Modelled	Actual	Comment	Reference
B	82	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	128	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	204	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	215	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
B	250	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	-2	SER	-	CLONING ARTIFACT	UNP Q81B18
C	-1	ASN	-	CLONING ARTIFACT	UNP Q81B18
C	0	ALA	-	CLONING ARTIFACT	UNP Q81B18
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	4	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	16	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	40	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	81	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	82	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	128	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	204	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	215	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
C	250	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	-2	SER	-	CLONING ARTIFACT	UNP Q81B18
D	-1	ASN	-	CLONING ARTIFACT	UNP Q81B18
D	0	ALA	-	CLONING ARTIFACT	UNP Q81B18
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	4	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	16	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	40	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	81	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	82	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	128	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	204	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	215	MSE	MET	MODIFIED RESIDUE	UNP Q81B18
D	250	MSE	MET	MODIFIED RESIDUE	UNP Q81B18

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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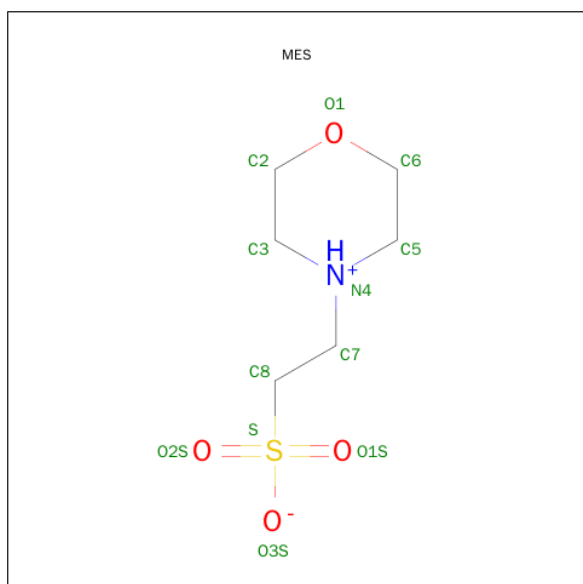
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

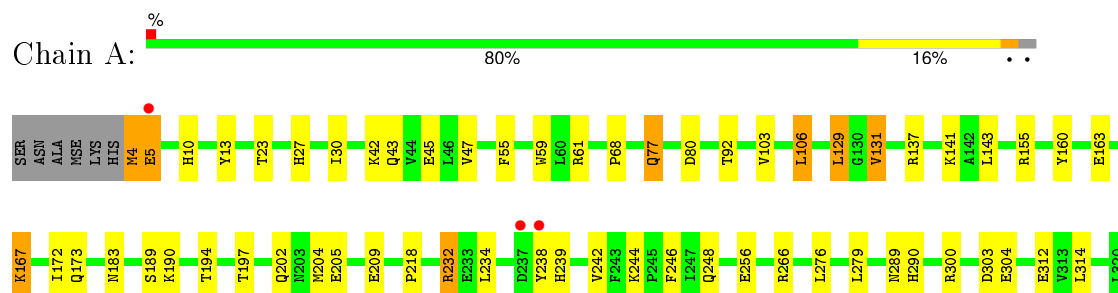
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	286	Total	O	0	0
			286	286		
5	B	301	Total	O	0	0
			301	301		
5	C	261	Total	O	0	0
			261	261		
5	D	196	Total	O	0	0
			196	196		

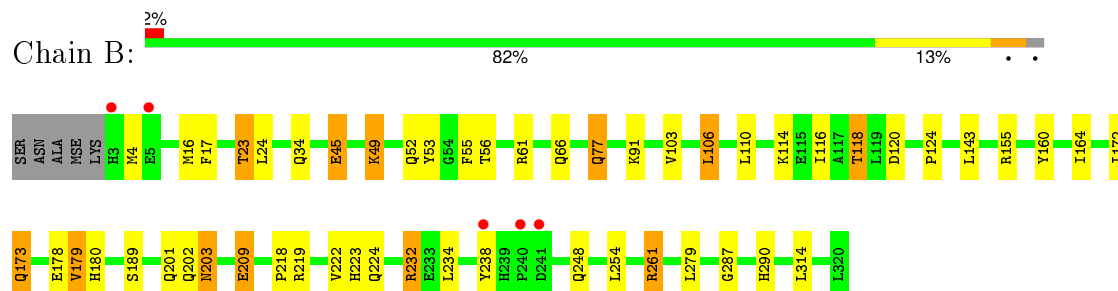
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

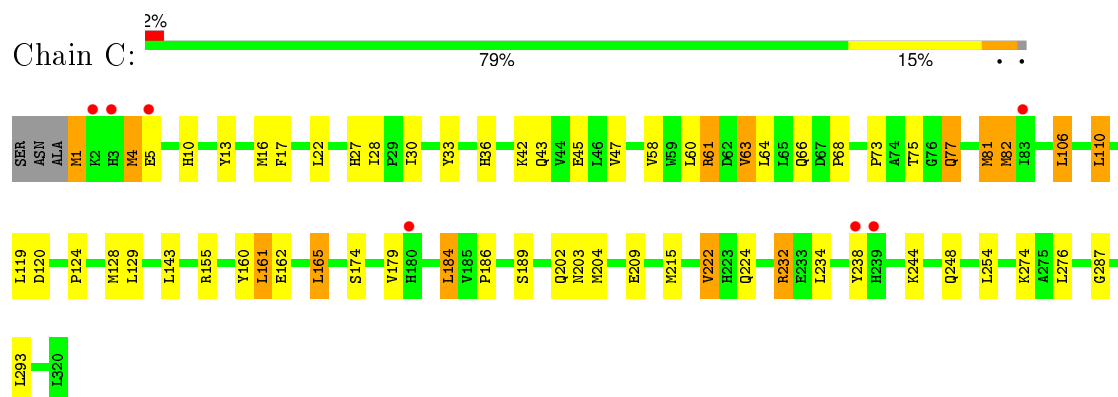
- Molecule 1: Luciferase-like monooxygenase



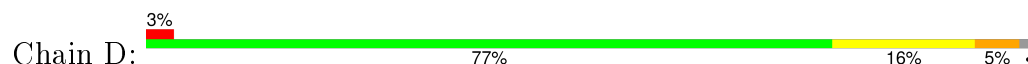
- Molecule 1: Luciferase-like monooxygenase

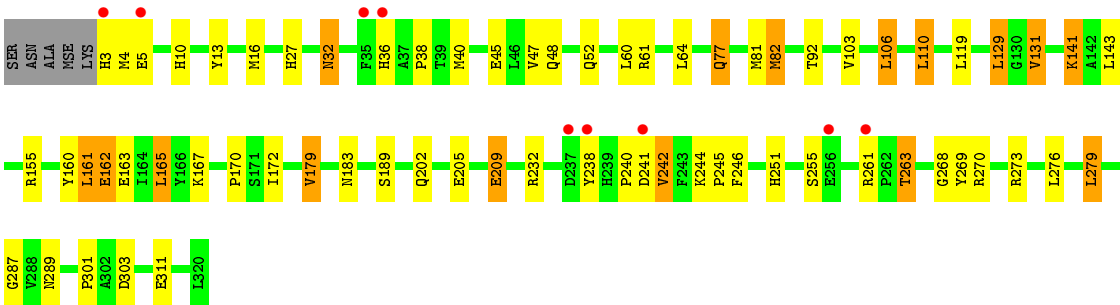


- Molecule 1: Luciferase-like monooxygenase



- Molecule 1: Luciferase-like monooxygenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	138.47Å 159.94Å 172.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.81 – 2.50 47.81 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.81-2.50) 99.5 (47.81-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.183 , 0.218 0.181 , 0.216	Depositor DCC
R_{free} test set	13183 reflections (11.14%)	DCC
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 131506 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12077	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, MES

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.58	0/2756	0.67	0/3719
1	B	0.58	0/2755	0.67	0/3718
1	C	0.57	0/2735	0.69	3/3689 (0.1%)
1	D	0.55	0/2762	0.62	0/3726
All	All	0.57	0/11008	0.66	3/14852 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	222	VAL	CB-CA-C	-6.00	100.01	111.40
1	C	61	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	C	61	ARG	NE-CZ-NH1	5.15	122.87	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2690	0	2592	48	0
1	B	2690	0	2592	45	0
1	C	2670	0	2587	51	0
1	D	2698	0	2602	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	60	0	0	0	0
2	B	30	0	0	1	0
2	C	55	0	0	0	0
2	D	50	0	0	0	0
3	A	24	0	24	6	0
3	B	12	0	12	0	0
3	C	12	0	12	0	0
3	D	12	0	12	0	0
4	A	12	0	16	0	0
4	B	6	0	8	2	0
4	C	6	0	8	2	0
4	D	6	0	8	0	0
5	A	286	0	0	5	0
5	B	301	0	0	9	0
5	C	261	0	0	3	0
5	D	196	0	0	8	0
All	All	12077	0	10473	193	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (193) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:MSE:HA	1:D:82:MSE:HE3	1.37	1.04
1:A:80:ASP:H	1:B:118:THR:HG21	1.27	0.99
3:A:2404:MES:H62	1:C:68:PRO:HA	1.48	0.95
1:D:38:PRO:HG2	1:D:40:MSE:HE2	1.51	0.93
1:B:155:ARG:HH21	1:B:202:GLN:HE22	1.18	0.88
1:C:155:ARG:HH11	1:C:202:GLN:HE22	1.20	0.87
1:B:173[A]:GLN:NE2	1:B:178[A]:GLU:OE2	2.07	0.87
1:D:40:MSE:SE	5:D:2641:HOH:O	2.42	0.87
1:B:61:ARG:H	1:B:77:GLN:HE22	1.20	0.86
1:C:61:ARG:H	1:C:77:GLN:HE22	1.23	0.86
1:B:23:THR:HG22	1:B:290:HIS:HD1	1.40	0.86
1:D:246:PHE:H	1:D:289:ASN:HD22	1.21	0.85
1:D:10:HIS:HD2	1:D:13:TYR:H	1.24	0.84
1:D:82:MSE:HA	1:D:82:MSE:CE	2.09	0.82
1:A:10:HIS:HD2	1:A:13:TYR:H	1.25	0.82
1:C:30:ILE:H	1:C:43:GLN:HE22	1.27	0.82
1:A:163[A]:GLU:OE2	1:A:167:LYS:HE2	1.80	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ILE:H	1:A:43:GLN:HE22	1.26	0.80
1:C:64:LEU:HD13	1:D:110:LEU:HD13	1.64	0.79
1:D:155:ARG:HE	1:D:202:GLN:HE22	1.28	0.79
1:C:1:MSE:N	1:C:4:MSE:HE3	1.97	0.79
1:A:246:PHE:H	1:A:289:ASN:HD22	1.31	0.78
1:C:215:MSE:SE	5:C:2714:HOH:O	2.50	0.78
1:A:42:LYS:HD2	1:A:45[B]:GLU:OE2	1.84	0.77
1:A:155:ARG:HH21	1:A:202:GLN:HE22	1.33	0.76
1:D:61:ARG:H	1:D:77:GLN:HE22	1.30	0.76
1:A:205[B]:GLU:HG3	1:A:238:TYR:OH	1.86	0.75
3:A:2404:MES:C6	1:C:68:PRO:HA	2.18	0.74
1:C:110:LEU:HD13	1:D:64:LEU:HD13	1.72	0.71
1:C:274:LYS:HZ1	4:C:2413:GOL:H2	1.53	0.71
1:C:1:MSE:H2	1:C:4:MSE:HE3	1.53	0.70
1:C:224:GLN:HE21	1:C:248:GLN:NE2	1.89	0.70
1:C:82:MSE:HE3	1:C:82:MSE:HA	1.73	0.70
1:C:204:MSE:HE3	1:C:234:LEU:CD1	2.22	0.70
1:D:38:PRO:HG2	1:D:40:MSE:CE	2.20	0.69
1:C:28:ILE:HD13	1:C:58:VAL:HG13	1.73	0.69
1:A:239:HIS:O	1:A:242:VAL:HG12	1.92	0.69
2:B:2421:SO4:O4	5:B:2654:HOH:O	2.09	0.69
1:C:215:MSE:CG	5:C:2714:HOH:O	2.42	0.68
1:B:23:THR:CG2	1:B:290:HIS:HD1	2.08	0.67
1:A:61:ARG:H	1:A:77:GLN:HE22	1.41	0.66
1:C:82:MSE:HE2	1:C:119:LEU:HD11	1.77	0.66
1:C:61:ARG:HD3	5:C:2460:HOH:O	1.95	0.66
1:C:10:HIS:HD2	1:C:13:TYR:H	1.45	0.65
1:A:239:HIS:HB2	1:A:242:VAL:HG11	1.77	0.65
1:A:10:HIS:CD2	1:A:13:TYR:H	2.12	0.64
1:C:42:LYS:HD2	1:C:45:GLU:OE1	1.97	0.64
1:C:27:HIS:CE1	1:C:61:ARG:HD2	2.33	0.64
1:A:42:LYS:CD	1:A:45[B]:GLU:OE2	2.45	0.64
1:C:204:MSE:HE3	1:C:234:LEU:HD12	1.78	0.63
1:D:61:ARG:CA	1:D:81:MSE:HE3	2.29	0.63
1:C:82:MSE:HE2	1:C:119:LEU:CD1	2.28	0.63
1:C:16:MSE:SE	1:C:128:MSE:HE3	2.48	0.62
1:D:129:LEU:HD22	1:D:131:VAL:HG23	1.82	0.62
1:B:261[A]:ARG:HE	1:B:261[A]:ARG:H	1.46	0.61
3:A:2404:MES:H22	1:C:73:PRO:HA	1.82	0.61
1:B:172:ILE:HB	1:B:179:VAL:HG13	1.81	0.61
1:A:256[A]:GLU:HG3	5:A:2772:HOH:O	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:MSE:CE	1:D:119:LEU:HD11	2.31	0.60
1:C:10:HIS:CD2	1:C:13:TYR:H	2.19	0.60
4:B:2412:GOL:H2	1:D:170:PRO:HA	1.83	0.60
1:C:60:LEU:O	1:C:81:MSE:HE1	2.02	0.60
1:B:232:ARG:NH2	1:B:287:GLY:O	2.34	0.60
1:A:239:HIS:HB2	1:A:242:VAL:CG1	2.31	0.60
1:B:222:VAL:CG2	5:B:2579:HOH:O	2.49	0.60
1:B:222:VAL:HG23	5:B:2579:HOH:O	2.02	0.60
1:A:61:ARG:HD3	5:A:2598:HOH:O	2.02	0.59
1:B:224:GLN:HE21	1:B:248:GLN:NE2	1.99	0.59
1:D:209[A]:GLU:HG2	1:D:238:TYR:CE2	2.38	0.59
1:B:23:THR:HG23	1:B:56:THR:OG1	2.03	0.58
1:A:204:MSE:HG2	1:A:234:LEU:HD12	1.87	0.56
1:D:232:ARG:NH2	1:D:287:GLY:O	2.38	0.56
1:A:303:ASP:OD2	1:A:304:GLU:N	2.40	0.55
1:D:269:TYR:CD1	1:D:279:LEU:HD11	2.42	0.55
1:A:59:TRP:HE1	1:A:290:HIS:HE1	1.56	0.54
1:C:232:ARG:NH2	1:C:287:GLY:O	2.40	0.54
1:D:82:MSE:HE1	1:D:119:LEU:CD1	2.37	0.54
1:A:163[B]:GLU:HB3	1:A:172:ILE:HD11	1.89	0.54
1:C:16:MSE:SE	1:C:17:PHE:CZ	3.11	0.54
1:C:66:GLN:NE2	1:D:183:ASN:HD22	2.06	0.54
1:D:263:THR:HG22	5:D:2509:HOH:O	2.08	0.53
1:C:204:MSE:HE3	1:C:234:LEU:HD11	1.90	0.53
1:A:27:HIS:HE1	1:A:61:ARG:HD2	1.74	0.53
1:B:49:LYS:HE2	5:B:2693:HOH:O	2.09	0.52
1:D:205:GLU:O	1:D:209[A]:GLU:HG3	2.10	0.52
1:C:209[A]:GLU:OE2	1:C:238:TYR:HB3	2.09	0.52
1:A:209[A]:GLU:HG2	1:A:238:TYR:CG	2.45	0.52
1:A:23:THR:OG1	1:A:290:HIS:HD2	1.93	0.52
1:C:274:LYS:NZ	4:C:2413:GOL:H2	2.24	0.52
1:B:106:LEU:HG	1:B:143:LEU:HD12	1.91	0.52
1:C:1:MSE:H3	1:C:4:MSE:HE3	1.73	0.52
1:D:82:MSE:HE2	1:D:119:LEU:HD11	1.92	0.51
1:B:201:GLN:NE2	5:B:2669:HOH:O	2.42	0.51
1:A:137:ARG:HG2	5:A:2753:HOH:O	2.11	0.51
1:B:201:GLN:HB3	5:B:2669:HOH:O	2.10	0.51
1:C:234:LEU:O	1:C:238:TYR:HD1	1.93	0.51
1:A:256[A]:GLU:HG2	1:A:300:ARG:HD2	1.91	0.51
3:A:2404:MES:H62	1:C:68:PRO:CA	2.31	0.51
1:A:246:PHE:H	1:A:289:ASN:ND2	2.05	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:ILE:HD13	1:C:58:VAL:CG1	2.40	0.50
1:B:23:THR:CG2	1:B:56:THR:OG1	2.60	0.50
1:A:27:HIS:CE1	1:A:61:ARG:HD2	2.46	0.49
4:B:2412:GOL:H11	1:D:167:LYS:HE2	1.94	0.49
1:D:61:ARG:C	1:D:81:MSE:HE3	2.32	0.49
1:A:4:MSE:HG2	1:A:5:GLU:N	2.28	0.49
1:A:106:LEU:HG	1:A:143:LEU:HD12	1.94	0.48
1:D:32:ASN:H	1:D:32:ASN:HD22	1.60	0.48
1:A:141[A]:LYS:HD3	5:B:2615:HOH:O	2.13	0.48
1:D:82:MSE:HE1	1:D:119:LEU:HD11	1.94	0.48
1:D:45:GLU:HG3	5:D:2615:HOH:O	2.12	0.48
1:A:183:ASN:HD22	1:B:66:GLN:NE2	2.10	0.48
1:C:27:HIS:HE1	1:C:61:ARG:HD2	1.75	0.48
1:D:10:HIS:CD2	1:D:13:TYR:H	2.16	0.48
1:B:173[A]:GLN:HE21	1:B:178[A]:GLU:CD	2.17	0.48
1:B:209:GLU:HB2	1:B:238:TYR:CD1	2.48	0.48
1:D:47:VAL:HG12	1:D:92:THR:HG21	1.95	0.48
1:B:114:LYS:O	1:B:118:THR:CG2	2.62	0.47
1:C:209[A]:GLU:HB2	1:C:238:TYR:CD2	2.49	0.47
1:A:68:PRO:HB3	3:A:2404:MES:H51	1.97	0.47
1:D:106:LEU:HG	1:D:143:LEU:HD12	1.96	0.47
1:D:103:VAL:HG11	1:D:106:LEU:HD22	1.96	0.47
1:B:173[A]:GLN:HE21	1:B:178[A]:GLU:CG	2.26	0.47
1:A:61:ARG:CD	5:A:2598:HOH:O	2.62	0.47
1:C:106:LEU:HG	1:C:143:LEU:HD12	1.96	0.47
1:C:63:VAL:HG22	1:C:75:THR:HG22	1.96	0.47
1:C:184:LEU:HD13	1:C:186:PRO:O	2.15	0.47
1:D:52[B]:GLN:HB3	5:D:2491:HOH:O	2.14	0.47
1:A:232[A]:ARG:NH1	1:A:244:LYS:O	2.48	0.47
1:B:34:GLN:NE2	5:B:2541:HOH:O	2.46	0.47
1:D:141:LYS:HD2	1:D:141:LYS:HA	1.53	0.46
1:D:61:ARG:HA	1:D:81:MSE:HE3	1.95	0.46
1:C:232:ARG:NH1	1:C:244:LYS:O	2.48	0.46
1:B:45[B]:GLU:O	1:B:49:LYS:HG3	2.15	0.46
1:D:3:HIS:HD2	1:D:242:VAL:HG11	1.81	0.46
1:D:161:LEU:O	1:D:165:LEU:HB2	2.16	0.46
1:B:4:MSE:H	1:B:4:MSE:SE	2.49	0.45
1:B:16:MSE:SE	1:B:17:PHE:CZ	3.19	0.45
1:B:52[C]:GLN:HG2	1:B:53:TYR:N	2.32	0.45
1:A:4:MSE:CE	1:A:5:GLU:H	2.29	0.45
1:A:55:PHE:CZ	1:A:314:LEU:HD21	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:PHE:N	1:A:289:ASN:HD22	2.08	0.44
1:C:161:LEU:O	1:C:165:LEU:HB2	2.18	0.44
1:D:163[A]:GLU:HG2	1:D:167:LYS:HD2	1.99	0.44
1:A:103:VAL:HG11	1:A:106:LEU:HD22	1.98	0.44
1:D:162:GLU:HG2	5:D:2592:HOH:O	2.17	0.44
1:C:120:ASP:O	1:C:124:PRO:HA	2.17	0.44
1:C:82:MSE:CE	1:C:82:MSE:HA	2.46	0.44
1:D:45:GLU:HA	1:D:48:GLN:HE21	1.83	0.43
1:D:172:ILE:HB	1:D:179:VAL:HG13	1.99	0.43
1:D:60:LEU:O	1:D:81:MSE:CE	2.67	0.43
1:D:255:SER:HB2	1:D:270:ARG:NH2	2.33	0.43
1:B:103:VAL:HG11	1:B:106:LEU:HD22	2.00	0.43
1:C:60:LEU:C	1:C:81:MSE:HE1	2.37	0.43
1:B:155:ARG:HD3	1:B:202:GLN:NE2	2.33	0.43
1:A:30:ILE:H	1:A:43:GLN:NE2	2.05	0.43
1:B:222:VAL:HG22	5:B:2579:HOH:O	2.17	0.43
1:B:223:HIS:HD2	1:B:223:HIS:O	2.02	0.43
1:C:33:TYR:HB2	1:C:36:HIS:O	2.19	0.43
1:A:59:TRP:HE1	1:A:290:HIS:CE1	2.34	0.43
1:B:218:PRO:O	1:B:219:ARG:HD3	2.19	0.42
3:A:2404:MES:C5	1:C:68:PRO:HA	2.49	0.42
1:D:36:HIS:HD2	5:D:2625:HOH:O	2.02	0.42
1:B:178[A]:GLU:CG	1:B:180:HIS:CD2	3.02	0.42
1:D:311:GLU:HG2	5:D:2643:HOH:O	2.19	0.42
1:B:114:LYS:O	1:B:118:THR:HG23	2.19	0.42
1:D:244:LYS:HA	1:D:245:PRO:HD3	1.88	0.42
1:A:218:PRO:HA	1:A:248:GLN:NE2	2.34	0.42
1:D:82:MSE:CE	1:D:82:MSE:CA	2.86	0.42
1:B:55:PHE:CZ	1:B:314:LEU:HD21	2.55	0.42
1:D:232:ARG:NH1	1:D:244:LYS:O	2.53	0.41
1:C:60:LEU:HB2	1:C:81:MSE:CE	2.50	0.41
1:A:129:LEU:HD22	1:A:131:VAL:HG23	2.01	0.41
1:A:47:VAL:HG12	1:A:92:THR:HG21	2.02	0.41
1:B:116:ILE:HA	1:B:116:ILE:HD13	1.95	0.41
1:A:209[A]:GLU:HG2	1:A:238:TYR:CD1	2.56	0.41
1:C:28:ILE:HD11	1:C:47:VAL:CG2	2.51	0.41
1:A:256[B]:GLU:HG3	5:A:2764:HOH:O	2.20	0.41
1:A:129:LEU:O	1:A:194:THR:HA	2.20	0.41
1:D:16:MSE:SE	5:D:2640:HOH:O	2.88	0.41
1:B:155:ARG:HD3	1:B:202:GLN:HE22	1.86	0.41
1:D:27:HIS:HE1	1:D:61:ARG:HG3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:PRO:O	1:D:241:ASP:HB2	2.21	0.41
1:B:203:ASN:HD22	1:B:203:ASN:C	2.24	0.41
1:A:30:ILE:N	1:A:43:GLN:HE22	2.06	0.40
1:B:234:LEU:O	1:B:238:TYR:HD2	2.04	0.40
1:A:4:MSE:HG2	1:A:5:GLU:HG3	2.03	0.40
1:B:110:LEU:HD13	1:B:164:ILE:HG12	2.03	0.40
1:D:251:HIS:O	1:D:268:GLY:HA3	2.22	0.40
1:B:120:ASP:O	1:B:124:PRO:HA	2.21	0.40
1:D:301:PRO:HB2	1:D:303[B]:ASP:OD1	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	329/323 (102%)	316 (96%)	13 (4%)	0	100	100
1	B	329/323 (102%)	319 (97%)	10 (3%)	0	100	100
1	C	327/323 (101%)	316 (97%)	11 (3%)	0	100	100
1	D	330/323 (102%)	320 (97%)	10 (3%)	0	100	100
All	All	1315/1292 (102%)	1271 (97%)	44 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	286/267 (107%)	266 (93%)	20 (7%)	19	34
1	B	286/267 (107%)	264 (92%)	22 (8%)	16	30
1	C	284/267 (106%)	258 (91%)	26 (9%)	11	21
1	D	287/267 (108%)	260 (91%)	27 (9%)	11	20
All	All	1143/1068 (107%)	1048 (92%)	95 (8%)	16	26

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	MSE
1	A	5	GLU
1	A	77	GLN
1	A	106	LEU
1	A	129	LEU
1	A	131	VAL
1	A	160	TYR
1	A	167	LYS
1	A	173[A]	GLN
1	A	173[B]	GLN
1	A	189[A]	SER
1	A	189[B]	SER
1	A	190	LYS
1	A	197	THR
1	A	232[A]	ARG
1	A	232[B]	ARG
1	A	266	ARG
1	A	276	LEU
1	A	279	LEU
1	A	312	GLU
1	B	23	THR
1	B	24	LEU
1	B	45[A]	GLU
1	B	45[B]	GLU
1	B	49	LYS
1	B	77	GLN
1	B	91	LYS
1	B	106	LEU
1	B	118	THR
1	B	160	TYR
1	B	173[A]	GLN
1	B	173[B]	GLN
1	B	179	VAL

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Mol	Chain	Res	Type
1	B	189[A]	SER
1	B	189[B]	SER
1	B	203	ASN
1	B	209	GLU
1	B	232	ARG
1	B	254	LEU
1	B	261[A]	ARG
1	B	261[B]	ARG
1	B	279	LEU
1	C	1	MSE
1	C	4	MSE
1	C	5	GLU
1	C	22	LEU
1	C	63	VAL
1	C	77	GLN
1	C	81	MSE
1	C	82	MSE
1	C	106	LEU
1	C	110	LEU
1	C	129	LEU
1	C	160	TYR
1	C	161	LEU
1	C	162	GLU
1	C	165	LEU
1	C	174	SER
1	C	179	VAL
1	C	184	LEU
1	C	189[A]	SER
1	C	189[B]	SER
1	C	203	ASN
1	C	222	VAL
1	C	232	ARG
1	C	254	LEU
1	C	276	LEU
1	C	293	LEU
1	D	4	MSE
1	D	5	GLU
1	D	32	ASN
1	D	77	GLN
1	D	82	MSE
1	D	106	LEU
1	D	110	LEU

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Mol	Chain	Res	Type
1	D	129	LEU
1	D	131	VAL
1	D	141	LYS
1	D	160	TYR
1	D	161	LEU
1	D	162	GLU
1	D	165	LEU
1	D	179	VAL
1	D	189[A]	SER
1	D	189[B]	SER
1	D	209[A]	GLU
1	D	209[B]	GLU
1	D	242	VAL
1	D	261[A]	ARG
1	D	261[B]	ARG
1	D	261[C]	ARG
1	D	263	THR
1	D	273	ARG
1	D	276	LEU
1	D	279	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	HIS
1	A	27	HIS
1	A	43	GLN
1	A	66	GLN
1	A	77	GLN
1	A	202	GLN
1	A	289	ASN
1	A	290	HIS
1	B	3	HIS
1	B	34	GLN
1	B	66	GLN
1	B	77	GLN
1	B	180	HIS
1	B	201	GLN
1	B	202	GLN
1	B	203	ASN
1	B	223	HIS
1	B	248	GLN

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Mol	Chain	Res	Type
1	C	10	HIS
1	C	27	HIS
1	C	43	GLN
1	C	66	GLN
1	C	77	GLN
1	C	202	GLN
1	C	203	ASN
1	C	248	GLN
1	D	3	HIS
1	D	10	HIS
1	D	27	HIS
1	D	32	ASN
1	D	48	GLN
1	D	77	GLN
1	D	147	HIS
1	D	202	GLN
1	D	239	HIS
1	D	289	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

49 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	MES	A	2403	-	11,12,12	0.72	0	14,16,16	2.76	5 (35%)
3	MES	A	2404	-	11,12,12	0.70	0	14,16,16	2.82	7 (50%)
4	GOL	A	2411	-	5,5,5	0.50	0	5,5,5	0.33	0
4	GOL	A	2415	-	5,5,5	0.32	0	5,5,5	0.61	0
2	SO4	A	2426	-	4,4,4	0.19	0	6,6,6	0.27	0
2	SO4	A	2427	-	4,4,4	0.16	0	6,6,6	0.12	0
2	SO4	A	2428	-	4,4,4	0.24	0	6,6,6	0.17	0
2	SO4	A	2429	-	4,4,4	0.15	0	6,6,6	0.22	0
2	SO4	A	2446	-	4,4,4	0.26	0	6,6,6	0.14	0
2	SO4	A	2447	-	4,4,4	0.15	0	6,6,6	0.40	0
2	SO4	A	2448	-	4,4,4	0.16	0	6,6,6	0.16	0
2	SO4	A	2449	-	4,4,4	0.21	0	6,6,6	0.15	0
2	SO4	A	2455	-	4,4,4	0.18	0	6,6,6	0.23	0
2	SO4	A	2530	-	4,4,4	0.12	0	6,6,6	0.22	0
2	SO4	A	2531	-	4,4,4	0.12	0	6,6,6	0.15	0
2	SO4	A	2532	-	4,4,4	0.13	0	6,6,6	0.20	0
3	MES	B	2405	-	11,12,12	0.76	0	14,16,16	2.44	4 (28%)
4	GOL	B	2412	-	5,5,5	0.40	0	5,5,5	0.36	0
2	SO4	B	2421	-	4,4,4	0.24	0	6,6,6	0.42	0
2	SO4	B	2433	-	4,4,4	0.18	0	6,6,6	0.17	0
2	SO4	B	2434	-	4,4,4	0.17	0	6,6,6	0.11	0
2	SO4	B	2442	-	4,4,4	0.19	0	6,6,6	0.14	0
2	SO4	B	2457	-	4,4,4	0.18	0	6,6,6	0.51	0
2	SO4	B	2458	-	4,4,4	0.15	0	6,6,6	0.13	0
3	MES	C	2401	-	11,12,12	0.77	0	14,16,16	2.45	4 (28%)
4	GOL	C	2413	-	5,5,5	0.35	0	5,5,5	0.39	0
2	SO4	C	2422	-	4,4,4	0.13	0	6,6,6	0.14	0
2	SO4	C	2436	-	4,4,4	0.15	0	6,6,6	0.21	0
2	SO4	C	2437	-	4,4,4	0.15	0	6,6,6	0.41	0
2	SO4	C	2438	-	4,4,4	0.16	0	6,6,6	0.12	0
2	SO4	C	2439	-	4,4,4	0.13	0	6,6,6	0.12	0
2	SO4	C	2444	-	4,4,4	0.16	0	6,6,6	0.25	0
2	SO4	C	2450	-	4,4,4	0.12	0	6,6,6	0.25	0
2	SO4	C	2452	-	4,4,4	0.19	0	6,6,6	0.13	0
2	SO4	C	2454	-	4,4,4	0.23	0	6,6,6	0.33	0
2	SO4	C	2456	-	4,4,4	0.20	0	6,6,6	0.19	0
2	SO4	C	2459	-	4,4,4	0.19	0	6,6,6	0.12	0
3	MES	D	2402	-	11,12,12	0.70	0	14,16,16	2.78	6 (42%)
4	GOL	D	2414	-	5,5,5	0.35	0	5,5,5	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	D	2423	-	4,4,4	0.08	0	6,6,6	0.27	0
2	SO4	D	2424	-	4,4,4	0.23	0	6,6,6	0.19	0
2	SO4	D	2425	-	4,4,4	0.14	0	6,6,6	0.12	0
2	SO4	D	2435	-	4,4,4	0.13	0	6,6,6	0.36	0
2	SO4	D	2440	-	4,4,4	0.16	0	6,6,6	0.24	0
2	SO4	D	2441	-	4,4,4	0.12	0	6,6,6	0.27	0
2	SO4	D	2443	-	4,4,4	0.21	0	6,6,6	0.30	0
2	SO4	D	2445	-	4,4,4	0.21	0	6,6,6	0.14	0
2	SO4	D	2451	-	4,4,4	0.15	0	6,6,6	0.18	0
2	SO4	D	2453	-	4,4,4	0.14	0	6,6,6	0.25	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MES	A	2403	-	-	0/6/14/14	0/1/1/1
3	MES	A	2404	-	-	0/6/14/14	0/1/1/1
4	GOL	A	2411	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2415	-	-	0/4/4/4	0/0/0/0
2	SO4	A	2426	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2427	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2428	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2429	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2446	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2447	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2448	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2449	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2455	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2530	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2531	-	-	0/0/0/0	0/0/0/0
2	SO4	A	2532	-	-	0/0/0/0	0/0/0/0
3	MES	B	2405	-	-	0/6/14/14	0/1/1/1
4	GOL	B	2412	-	-	0/4/4/4	0/0/0/0
2	SO4	B	2421	-	-	0/0/0/0	0/0/0/0
2	SO4	B	2433	-	-	0/0/0/0	0/0/0/0
2	SO4	B	2434	-	-	0/0/0/0	0/0/0/0
2	SO4	B	2442	-	-	0/0/0/0	0/0/0/0
2	SO4	B	2457	-	-	0/0/0/0	0/0/0/0
2	SO4	B	2458	-	-	0/0/0/0	0/0/0/0
3	MES	C	2401	-	-	0/6/14/14	0/1/1/1
4	GOL	C	2413	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	C	2422	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2436	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2437	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2438	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2439	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2444	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2450	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2452	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2454	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2456	-	-	0/0/0/0	0/0/0/0
2	SO4	C	2459	-	-	0/0/0/0	0/0/0/0
3	MES	D	2402	-	-	0/6/14/14	0/1/1/1
4	GOL	D	2414	-	-	0/4/4/4	0/0/0/0
2	SO4	D	2423	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2424	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2425	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2435	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2440	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2441	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2443	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2445	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2451	-	-	0/0/0/0	0/0/0/0
2	SO4	D	2453	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2404	MES	C6-C5-N4	-4.62	103.12	110.12
3	D	2402	MES	C6-C5-N4	-4.58	103.18	110.12
3	A	2404	MES	C2-C3-N4	-4.53	103.27	110.12
3	D	2402	MES	C2-C3-N4	-3.67	104.57	110.12
3	A	2404	MES	O1-C6-C5	-2.50	106.11	111.84
3	C	2401	MES	C7-N4-C3	2.15	116.77	111.27
3	A	2403	MES	C7-N4-C5	2.29	117.14	111.27
3	A	2403	MES	C7-N4-C3	2.31	117.18	111.27
3	B	2405	MES	C7-N4-C5	2.48	117.63	111.27
3	B	2405	MES	C7-N4-C3	2.65	118.06	111.27
3	C	2401	MES	C7-N4-C5	2.66	118.10	111.27
3	A	2404	MES	O2S-S-C8	2.71	109.22	106.91
3	C	2401	MES	O1S-S-C8	2.76	109.26	106.91
3	A	2403	MES	C2-C3-N4	2.91	114.53	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2402	MES	C7-N4-C3	3.19	119.45	111.27
3	D	2402	MES	C7-N4-C5	3.46	120.14	111.27
3	B	2405	MES	O1S-S-C8	3.48	109.87	106.91
3	A	2404	MES	C7-N4-C5	3.67	120.67	111.27
3	A	2404	MES	C7-N4-C3	3.77	120.94	111.27
3	D	2402	MES	O1S-S-C8	3.78	110.13	106.91
3	A	2404	MES	C5-N4-C3	4.38	118.38	108.90
3	A	2403	MES	O2S-S-C8	4.44	110.69	106.91
3	D	2402	MES	C5-N4-C3	5.31	120.39	108.90
3	B	2405	MES	C5-N4-C3	7.11	124.30	108.90
3	C	2401	MES	C5-N4-C3	7.50	125.13	108.90
3	A	2403	MES	C5-N4-C3	7.75	125.68	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2404	MES	6	0
4	B	2412	GOL	2	0
2	B	2421	SO4	1	0
4	C	2413	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	308/323 (95%)	-0.44	3 (0%) 84 86	20, 29, 42, 65	0
1	B	309/323 (95%)	-0.45	5 (1%) 74 78	20, 28, 44, 81	0
1	C	310/323 (95%)	-0.34	7 (2%) 64 67	22, 31, 46, 100	0
1	D	309/323 (95%)	-0.17	9 (2%) 55 60	22, 37, 54, 83	0
All	All	1236/1292 (95%)	-0.35	24 (1%) 70 73	20, 31, 49, 100	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3	HIS	5.6
1	B	3	HIS	5.4
1	C	3	HIS	5.1
1	B	5	GLU	3.7
1	C	238	TYR	3.5
1	A	238	TYR	3.4
1	B	238	TYR	3.4
1	C	5	GLU	3.3
1	D	36	HIS	3.3
1	D	238	TYR	3.2
1	C	2	LYS	2.9
1	B	241	ASP	2.8
1	C	180	HIS	2.8
1	A	5	GLU	2.7
1	D	35	PHE	2.5
1	D	5	GLU	2.5
1	C	239	HIS	2.4
1	C	83	ILE	2.4
1	D	256[A]	GLU	2.2
1	B	240	PRO	2.2
1	D	241	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	237	ASP	2.1
1	A	237	ASP	2.1
1	D	261[A]	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	A	2532	5/5	0.84	0.34	9.40	113,113,113,114	0
2	SO4	D	2440	5/5	0.94	0.49	9.29	96,96,96,97	0
4	GOL	B	2412	6/6	0.87	0.31	6.33	51,53,54,56	0
2	SO4	A	2428	5/5	0.95	0.28	6.08	98,98,98,98	0
3	MES	A	2404	12/12	0.78	0.26	5.68	63,70,82,82	0
4	GOL	A	2415	6/6	0.86	0.23	4.45	69,69,70,71	0
4	GOL	D	2414	6/6	0.87	0.24	4.14	52,54,54,54	0
2	SO4	A	2427	5/5	0.93	0.22	3.55	78,78,78,78	0
2	SO4	A	2531	5/5	0.90	0.22	2.73	95,95,96,96	0
2	SO4	C	2444	5/5	0.89	0.18	1.88	100,100,101,101	0
2	SO4	D	2445	5/5	0.94	0.37	1.80	116,116,116,116	0
2	SO4	C	2422	5/5	0.92	0.30	1.71	73,74,74,74	1
2	SO4	D	2451	5/5	0.98	0.13	1.63	59,61,61,62	0
2	SO4	D	2441	5/5	0.97	0.23	1.55	81,81,82,82	0
2	SO4	A	2530	5/5	0.86	0.29	1.24	92,92,93,93	0
2	SO4	B	2457	5/5	0.94	0.23	1.12	67,68,68,70	0
2	SO4	A	2449	5/5	0.90	0.20	1.09	105,105,105,106	0
3	MES	B	2405	12/12	0.95	0.14	0.59	58,62,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MES	A	2403	12/12	0.97	0.13	0.43	51,55,57,58	0
2	SO4	B	2458	5/5	0.87	0.19	0.31	101,102,102,103	0
2	SO4	C	2437	5/5	0.94	0.19	0.15	74,75,75,75	0
2	SO4	C	2454	5/5	0.93	0.13	0.14	65,67,68,68	0
2	SO4	A	2426	5/5	0.97	0.19	0.08	74,74,75,75	0
2	SO4	A	2447	5/5	0.95	0.16	-0.02	73,73,73,73	0
2	SO4	D	2424	5/5	0.96	0.18	-0.06	77,78,78,78	0
3	MES	C	2401	12/12	0.97	0.14	-0.09	45,53,58,58	0
2	SO4	B	2421	5/5	0.99	0.11	-0.51	51,51,52,53	0
2	SO4	D	2435	5/5	0.94	0.13	-0.53	64,65,66,66	0
2	SO4	A	2429	5/5	0.95	0.10	-0.64	73,73,74,74	0
2	SO4	D	2443	5/5	0.95	0.11	-0.79	71,71,72,72	0
2	SO4	D	2423	5/5	0.97	0.09	-1.48	53,54,55,56	0
2	SO4	D	2425	5/5	0.86	0.30	-	99,99,99,99	1
2	SO4	C	2436	5/5	0.90	0.30	-	94,94,94,94	0
2	SO4	B	2442	5/5	0.91	0.26	-	85,85,86,86	0
2	SO4	C	2438	5/5	0.94	0.22	-	89,90,90,90	0
2	SO4	C	2459	5/5	0.84	0.45	-	125,125,126,126	0
2	SO4	A	2455	5/5	0.80	0.34	-	112,112,113,113	0
2	SO4	A	2448	5/5	0.79	0.27	-	102,102,102,103	0
2	SO4	B	2434	5/5	0.94	0.34	-	92,93,93,93	0
2	SO4	B	2433	5/5	0.92	0.21	-	99,99,100,100	0
2	SO4	D	2453	5/5	0.91	0.27	-	98,98,99,99	0
2	SO4	C	2450	5/5	0.93	0.20	-	74,75,76,76	0
2	SO4	C	2456	5/5	0.93	0.26	-	98,98,98,99	0
4	GOL	C	2413	6/6	0.86	0.16	-	68,69,69,70	0
2	SO4	C	2452	5/5	0.91	0.36	-	122,122,122,122	0
3	MES	D	2402	12/12	0.73	0.31	-	111,112,115,115	0
2	SO4	C	2439	5/5	0.93	0.28	-	97,98,98,98	0
4	GOL	A	2411	6/6	0.91	0.27	-	52,55,56,57	0
2	SO4	A	2446	5/5	0.96	0.17	-	72,73,74,74	0

6.5 Other polymers

There are no such residues in this entry.